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NEWS EXPRESS April 4 CURRENT WINDOWS VERSION IS V6.01a, CURRENT MACINTOSH VERSION IS V6.0b(ENG) AND V6.0Jb(JP),

AND CURRENT DISCOVER FILE IS DATED 01 APRIL 2003 NEWS HOURS STN Operating Hours Plus Help Desk Availability

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=> file reg COST IN U.S. DOLLARS

SINCE FILE TOTAL ENTRY SESSION

0.21

0.21

FULL ESTIMATED COST

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STRUCTURE FILE UPDATES: 22 MAY 2003 HIGHEST RN 519137-84-9 DICTIONARY FILE UPDATES: 22 MAY 2003 HIGHEST RN 519137-84-9

TSCA INFORMATION NOW CURRENT THROUGH JANUARY 6, 2003

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Crossover limits have been increased. See HELP CROSSOVER for details.

Experimental and calculated property data are now available. See HELP PROPERTIES for more information. See STNote 27, Searching Properties in the CAS Registry File, for complete details: http://www.cas.org/ONLINE/STN/STNOTES/stnotes27.pdf

Uploading 10004867.str

L1 STRUCTURE UPLOADED

=> d l1 L1 HAS NO ANSWERS L1 STR

G1 Cb, Hy, Ak

Structure attributes must be viewed using STN Express query preparation.

=> s l1 ful

FULL SEARCH INITIATED 15:34:57 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 3443 TO ITERATE

100.0% PROCESSED 3443 ITERATIONS

13 ANSWERS

SEARCH TIME: 00.00.01

L2 13 SEA SSS FUL L1

=> file caplus

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ENTRY 148.15

SESSION 148.36

FULL ESTIMATED COST

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FILE COVERS 1907 - 23 May 2003 VOL 138 ISS 22 FILE LAST UPDATED: 22 May 2003 (20030522/ED)

This file contains CAS Registry Numbers for easy and accurate substance identification.

=> s 12

L3

1 L2

=> d l3 1- ibib abs hitstr YOU HAVE REQUESTED DATA FROM 1 ANSWERS - CONTINUE? Y/(N):y

ANSWER 1 OF 1 CAPLUS COPYRIGHT 2003 ACS 2002:594821 CAPLUS ACCESSION NUMBER:

DOCUMENT NUMBER: 137:154856

Preparation of N-indanyl sulfonamides as potassium TITLE: applicant's

channel inhibitors

Beaudoin, Serge; Reed, Aimee D.; Gross, Michael INVENTOR(S):

PATENT ASSIGNEE(S): Icagen Incorporated, USA PCT Int. Appl., 72 pp. SOURCE:

CODEN: PIXXD2

DOCUMENT TYPE:

Patent English

LANGUAGE: FAMILY ACC. NUM. COUNT:

PATENT INFORMATION:

PATENT NO. KIND DATE APPLICATION NO. DATE ____ WO 2002060874 A1 20020808 WO 2001-US48601 20011219 WO 2002060874 C1 20030220 W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NO, NZ, OM, PH, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TN, TR, TT, TZ, UA, UG, UZ, VN, YU, ZA, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AT, BE, CH, CY, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, → US 2002161011 A1 20021031 US 2001-4867 20011207 PRIORITY APPLN. INFO.: US 2000-256926P P 20001221 A 20011207 US 2001-4867

OTHER SOURCE(S):

MARPAT 137:154856

GI

* STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY - AVAILABLE VIA OFFLINE PRINT *

AB The title compds. [I; A, B, D = C, N, N(O) (wherein at least one of A, B, and D is a substituted C atom and at most only one of A, B, and D is N(O)); E = H, alkyl; G = H, alkyl; or E and G taken together form a bond (site of unsatn.); R1 = H, alkyl, aryl, etc.; R2 = alkyl, aryl, heterocyclyl; R3 = H, alkyl, aryl, etc.; R4 = alkyl, aryl, heteroaryl, etc.; R5, R6 = H, F, alkyl; or R5 and R6 taken together, along with the carbom atom to which they are both attached, form a 3-7 membered carbocyclic or heterocyclic ring; R7 = H, alkyl, OH, etc.; n = 1-3], useful as potassium channel inhibitors and esp. useful for the treatment of cardiac arrhythmias and cell proliferative disorders, were prepd. Thus, amidation of the amine II (prepn. given) with hydrocinnamoyl chloride in the presence of Et3N in THF afforded 21% III which showed 46% inhibition of Kv1.5 at 0.1 .mu.M.

445402-76-6P 445402-81-3P 445402-82-4P IT 445402-85-7P 445402-86-8P

> RL: PAC (Pharmacological activity); RCT (Reactant); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); RACT (Reactant or reagent); USES (Uses)

(prepn. of N-indanyl sulfonamides as potassium channel inhibitors)

RN 445402-76-6 CAPLUS

CN 3-Pyridinecarboxylic acid, 2-[[3-[[(4-ethylphenyl)sulfonyl]amino]-2,3dihydro-1H-inden-5-yl]amino]- (9CI) (CA INDEX NAME)

RN 445402-81-3 CAPLUS

CN 3-Pyridinecarboxylic acid, 2-[[3-[[(4-fluorophenyl)sulfonyl]amino]-2,3-dihydro-1H-inden-5-yl]amino]-, ethyl ester (9CI) (CA INDEX NAME)

RN 445402-82-4 CAPLUS

CN 3-Pyridinecarboxylic acid, 2-[[3-[[(4-fluorophenyl)sulfonyl]amino]-2,3-dihydro-1H-inden-5-yl]amino]- (9CI) (CA INDEX NAME)

RN 445402-85-7 CAPLUS

CN Benzenesulfonamide, N-[2,3-dihydro-6-[(3-nitro-2-pyridinyl)amino]-1H-inden-1-yl]-4-ethyl- (9CI) (CA INDEX NAME)

RN 445402-86-8 CAPLUS

CN Benzenesulfonamide, N-[6-[(3-amino-2-pyridinyl)amino]-2,3-dihydro-1H-inden-1-yl]-4-ethyl- (9CI) (CA INDEX NAME)

IT 445402-77-7P 445402-78-8P 445402-79-9P 445402-80-2P 445402-83-5P 445402-84-6P 445402-87-9P 445402-88-0P

RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(prepn. of N-indanyl sulfonamides as potassium channel inhibitors)

RN 445402-77-7 CAPLUS

CN 3-Pyridinecarboxamide, 2-[[3-[[(4-ethylphenyl)sulfonyl]amino]-2,3-dihydro-1H-inden-5-yl]amino]-N-2-propenyl- (9CI) (CA INDEX NAME)

RN 445402-78-8 CAPLUS

CN 3-Pyridinecarboxamide, 2-[[3-[[(4-ethylphenyl)sulfonyl]amino]-2,3-dihydro-1H-inden-5-yl]amino]-N-(2-furanylmethyl)- (9CI) (CA INDEX NAME)

RN 445402-79-9 CAPLUS

CN 3-Pyridinecarboxamide, 2-[[3-[[(4-ethylphenyl)sulfonyl]amino]-2,3-dihydro-1H-inden-5-yl]amino]-N-2-propynyl- (9CI) (CA INDEX NAME)

RN 445402-80-2 CAPLUS

CN 3-Pyridinecarboxamide, 2-[[3-[[(4-ethylphenyl)sulfonyl]amino]-2,3-dihydro-1H-inden-5-yllamino]-N-propyl- (9CI) (CA INDEX NAME)

RN 445402-83-5 CAPLUS

CN 3-Pyridinecarboxamide, 2-[[3-[[(4-fluorophenyl)sulfonyl]amino]-2,3-dihydro-1H-inden-5-yl]amino]-N-(2-phenylethyl)- (9CI) (CA INDEX NAME)

RN 445402-84-6 CAPLUS

CN 3-Pyridinecarboxamide, 2-[[3-[[(4-fluorophenyl)sulfonyl]amino]-2,3-dihydro-1H-inden-5-yl]amino]-N-2-propenyl- (9CI) (CA INDEX NAME)

RN 445402-87-9 CAPLUS

CN 3-Pyridinecarboxamide, 2-[[3-[[(4-ethylphenyl)sulfonyl]amino]-2,3-dihydro-1H-inden-5-yl]amino]-N-(2-phenylethyl)- (9CI) (CA INDEX NAME)

RN 445402-88-0 CAPLUS

CN 3-Pyridinecarboxamide, 2-[[3-[[(4-fluorophenyl)sulfonyl]amino]-2,3-dihydro-1H-inden-5-yl]amino]-N-(2-thienylmethyl) (9CI) (CA INDEX NAME)

REFERENCE COUNT: 6 THERE ARE 6 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

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(FILE 'HOME' ENTERED AT 15:33:42 ON 23 MAY 2003)

FILE 'REGISTRY' ENTERED AT 15:34:27 ON 23 MAY 2003

L1 STRUCTURE UPLOADED

L2 13 S L1 FUL

FILE 'CAPLUS' ENTERED AT 15:35:05 ON 23 MAY 2003 L3 1 S L2

=> log y

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SINCE FILE TOTAL
ENTRY SESSION
FULL ESTIMATED COST

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)

SINCE FILE TOTAL

ENTRY SESSION

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